

Connecting via Winsock to STN

Welcome to STN International! Enter x:x

LOGINID:SSPTAKAB1626

PASSWORD:

TERMINAL (ENTER 1, 2, 3, OR ?):2

* * * * * Welcome to STN International * * * * *

NEWS	1		Web Page for STN Seminar Schedule - N. America
NEWS	2	DEC 01	ChemPort single article sales feature unavailable
NEWS	3	FEB 02	Simultaneous left and right truncation (SLART) added for CERAB, COMPUAB, ELCOM, and SOLIDSTATE
NEWS	4	FEB 02	GENBANK enhanced with SET PLURALS and SET SPELLING
NEWS	5	FEB 06	Patent sequence location (PSL) data added to USGENE
NEWS	6	FEB 10	COMPENDEX reloaded and enhanced
NEWS	7	FEB 11	WTEXTILES reloaded and enhanced
NEWS	8	FEB 19	New patent-examiner citations in 300,000 CA/CAPLUS patent records provide insights into related prior art
NEWS	9	FEB 19	Increase the precision of your patent queries -- use terms from the IPC Thesaurus, Version 2009.01
NEWS	10	FEB 23	Several formats for image display and print options discontinued in USPATFULL and USPAT2
NEWS	11	FEB 23	MEDLINE now offers more precise author group fields and 2009 MeSH terms
NEWS	12	FEB 23	TOXCENTER updates mirror those of MEDLINE - more precise author group fields and 2009 MeSH terms
NEWS	13	FEB 23	Three million new patent records blast AEROSPACE into STN patent clusters
NEWS	14	FEB 25	USGENE enhanced with patent family and legal status display data from INPADOCDB
NEWS	15	MAR 06	INPADOCDB and INPAFAMDB enhanced with new display formats
NEWS	16	MAR 11	EPFULL backfile enhanced with additional full-text applications and grants
NEWS	17	MAR 11	ESBIOBASE reloaded and enhanced
NEWS	18	MAR 20	CAS databases on STN enhanced with new super role for nanomaterial substances
NEWS	19	MAR 23	CA/CAPLUS enhanced with more than 250,000 patent equivalents from China
NEWS	20	MAR 30	IMSPATENTS reloaded and enhanced
NEWS	21	APR 03	CAS coverage of exemplified prophetic substances enhanced
NEWS	22	APR 07	STN is raising the limits on saved answers
NEWS	23	APR 24	CA/CAPLUS now has more comprehensive patent assignee information
NEWS	24	APR 26	USPATFULL and USPAT2 enhanced with patent assignment/reassignment information
NEWS	25	APR 28	CAS patent authority coverage expanded
NEWS	26	APR 28	ENCOMPLIT/ENCOMPLIT2 search fields enhanced
NEWS	27	APR 28	Limits doubled for structure searching in CAS

REGISTRY

NEWS EXPRESS JUNE 27 08 CURRENT WINDOWS VERSION IS V8.3,
AND CURRENT DISCOVER FILE IS DATED 23 JUNE 2008.

NEWS HOURS STN Operating Hours Plus Help Desk Availability
NEWS LOGIN Welcome Banner and News Items

Enter NEWS followed by the item number or name to see news on that specific topic.

All use of STN is subject to the provisions of the STN customer agreement. This agreement limits use to scientific research. Use for software development or design, implementation of commercial gateways, or use of CAS and STN data in the building of commercial products is prohibited and may result in loss of user privileges and other penalties.

* * * * * STN Columbus * * * * *

FILE 'HOME' ENTERED AT 13:39:19 ON 03 MAY 2009

=> file reg

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	0.22	0.22

FILE 'REGISTRY' ENTERED AT 13:39:30 ON 03 MAY 2009

USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.

PLEASE SEE "HELP USAGETERMS" FOR DETAILS.

COPYRIGHT (C) 2009 American Chemical Society (ACS)

Property values tagged with IC are from the ZIC/VINITI data file provided by InfoChem.

STRUCTURE FILE UPDATES: 1 MAY 2009 HIGHEST RN 1141923-26-3
DICTIONARY FILE UPDATES: 1 MAY 2009 HIGHEST RN 1141923-26-3

New CAS Information Use Policies, enter HELP USAGETERMS for details.

TSCA INFORMATION NOW CURRENT THROUGH January 9, 2009.

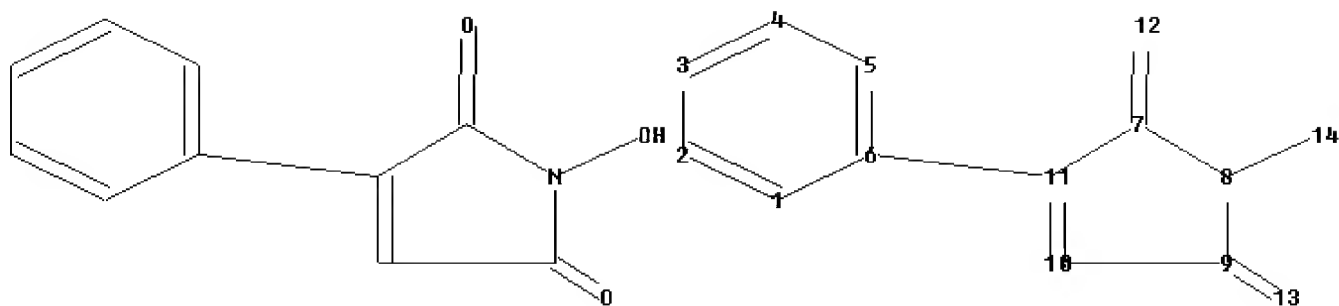
Please note that search-term pricing does apply when conducting SmartSELECT searches.

REGISTRY includes numerically searchable data for experimental and predicted properties as well as tags indicating availability of experimental property data in the original document. For information on property searching in REGISTRY, refer to:

<http://www.cas.org/support/stngen/stndoc/properties.html>

=>

Uploading C:\Program Files\STNEXP\Queries\10580230 unity.str



```

chain nodes :
12 13 14
ring nodes :
1 2 3 4 5 6 7 8 9 10 11
chain bonds :
6-11 7-12 8-14 9-13
ring bonds :
1-2 1-6 2-3 3-4 4-5 5-6 7-8 7-11 8-9 9-10 10-11
exact/norm bonds :
7-8 7-12 8-9 8-14 9-13
exact bonds :
6-11 7-11 9-10 10-11
normalized bonds :
1-2 1-6 2-3 3-4 4-5 5-6
isolated ring systems :
containing 7 :

```

```

Match level :
1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 10:Atom
11:Atom 12:CLASS 13:CLASS 14:CLASS

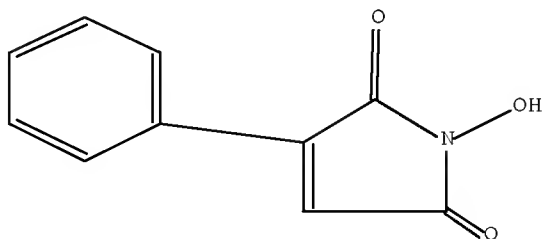
```

L1 STRUCTURE UPLOADED

=> d L1

L1 HAS NO ANSWERS

L1 STR



Structure attributes must be viewed using STN Express query preparation.

=> file caplus

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	0.48	0.70

FILE 'CAPLUS' ENTERED AT 13:39:46 ON 03 MAY 2009
 USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.
 PLEASE SEE "HELP USAGETERMS" FOR DETAILS.
 COPYRIGHT (C) 2009 AMERICAN CHEMICAL SOCIETY (ACS)

Copyright of the articles to which records in this database refer is held by the publishers listed in the PUBLISHER (PB) field (available for records published or updated in Chemical Abstracts after December 26, 1996), unless otherwise indicated in the original publications. The CA Lexicon is the copyrighted intellectual property of the American Chemical Society and is provided to assist you in searching databases on STN. Any dissemination, distribution, copying, or storing of this information, without the prior written consent of CAS, is strictly prohibited.

FILE COVERS 1907 - 3 May 2009 VOL 150 ISS 19
 FILE LAST UPDATED: 1 May 2009 (20090501/ED)

Caplus now includes complete International Patent Classification (IPC) reclassification data for the third quarter of 2008.

CAS Information Use Policies apply and are available at:

<http://www.cas.org/legal/infopolicy.html>

This file contains CAS Registry Numbers for easy and accurate substance identification.

=> s L1 SSS full

REGISTRY INITIATED

Substance data SEARCH and crossover from CAS REGISTRY in progress...
 Use DISPLAY HITSTR (or FHITSTR) to directly view retrieved structures.

FULL SEARCH INITIATED 13:39:51 FILE 'REGISTRY'
 FULL SCREEN SEARCH COMPLETED - 340 TO ITERATE

100.0% PROCESSED 340 ITERATIONS 25 ANSWERS
 SEARCH TIME: 00.00.01

L2 25 SEA SSS FUL L1

L3 17 L2

=> s L3 AND PY<=2003

24035378 PY<=2003

L4 3 L3 AND PY<=2003

=> d ibib abs hitstr 1-

YOU HAVE REQUESTED DATA FROM 3 ANSWERS - CONTINUE? Y/(N):y

L4 ANSWER 1 OF 3 CAPLUS COPYRIGHT 2009 ACS on STN

ACCESSION NUMBER: 2002:448755 CAPLUS Full-text

DOCUMENT NUMBER: 137:277867

TITLE: The himanimides, new bioactive compounds from *Serpula himantoides* (Fr.) Karst

AUTHOR(S): Aqueveque, Pedro; Anke, Timm; Sterner, Olov

CORPORATE SOURCE: Universidad de Concepcion, Concepcion, 3, Chile

SOURCE: Zeitschrift fuer Naturforschung, C: Journal of Biosciences (2002), 57(3/4), 257-262
CODEN: ZNCBDA; ISSN: 0939-5075

PUBLISHER: Verlag der Zeitschrift fuer Naturforschung

DOCUMENT TYPE: Journal

LANGUAGE: English

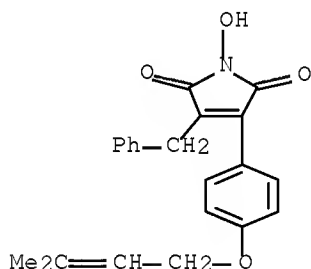
AB In a screening of basidiomycete cultures from Chile for the production of antibiotics the authors identified a *Serpula himantoides* strain as a producer of metabolites inhibiting the growth of bacteria and fungi. Bioactivity guided purification resulted in the isolation of 4 new antibiotics. Their structures were elucidated by spectroscopic methods. All 4 compds. are succinimide and maleimide derivs., of which 2 are N-hydroxylated.

IT 464189-92-2P, Himanimide C

RL: BPN (Biosynthetic preparation); BSU (Biological study, unclassified);
PRP (Properties); PUR (Purification or recovery); BIOL (Biological study);
PREP (Preparation)
(himanimide antibiotics from *Serpula himantoides* fermentation)

RN 464189-92-2 CAPLUS

CN 1H-Pyrrole-2,5-dione, 1-hydroxy-3-[4-[(3-methyl-2-buten-1-yl)oxy]phenyl]-4-(phenylmethyl)- (CA INDEX NAME)



REFERENCE COUNT: 8 THERE ARE 8 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L4 ANSWER 2 OF 3 CAPLUS COPYRIGHT 2009 ACS on STN

ACCESSION NUMBER: 1980:585380 CAPLUS Full-text

DOCUMENT NUMBER: 93:185380

ORIGINAL REFERENCE NO.: 93:29531a, 29534a

TITLE: Effects of solvents, N-substituents and acids on the photocyclization and the fluorescence behavior of diphenylmaleimides

AUTHOR(S): Ichimura, Kunihiro; Watanabe, Shoji; Kusakawa, Koichi; Ochi, Hideo

CORPORATE SOURCE: Res. Inst. Polym. Text., Ibaraki, 305, Japan

SOURCE: Nippon Kagaku Kaishi (1980), (6), 837-45

CODEN: NKAKB8; ISSN: 0369-4577

DOCUMENT TYPE: Journal

LANGUAGE: Japanese

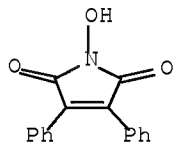
AB The fluorescence behavior of diphenylmaleimides shows the following characteristic features: (1) the Stokes shift correlates linearly with the solvent polarity parameter, ET, (2) the emission is quenched intramol. when the imido N is attached to atoms having n- or π -electrons, and (3) the quenching is observed in strongly acidic solns. The photocyclization of the imides to yield phenanthrenes and 9,10-dihydrophenanthrene-9,10-dicarboximides is influenced by the dual effect of acids; at pH about 3, the yield of the dihydrophenanthrenes increases with a decrease in pH without a change in the reaction rate, whereas in more strongly acidic solns. (H0 .apprx.2) the photoreactivity is reduced in proportion to the fluorescence quenching. The cyclization mechanism is discussed.

IT 75255-87-7

RL: RCT (Reactant); RACT (Reactant or reagent)
(fluorescence and photocyclization of)

RN 75255-87-7 CAPLUS

CN 1H-Pyrrole-2,5-dione, 1-hydroxy-3,4-diphenyl- (CA INDEX NAME)



L4 ANSWER 3 OF 3 CAPLUS COPYRIGHT 2009 ACS on STN

ACCESSION NUMBER: 1943:3617 CAPLUS Full-text

DOCUMENT NUMBER: 37:3617

ORIGINAL REFERENCE NO.: 37:641i,642a-c

TITLE: p-Bromophenylhydroxymaleic imide

AUTHOR(S): Skinner, Glenn S.; Coghlan, C. A.; Berlin, A. S.

SOURCE: Journal of the American Chemical Society (1942
, 64, 2600-1

CODEN: JACSAT; ISSN: 0002-7863

DOCUMENT TYPE: Journal

LANGUAGE: Unavailable

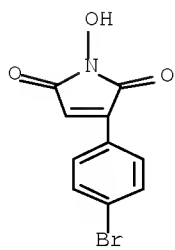
AB PhC(CN):C(OH)CO₂Et (21.7 g.) in 80 cc. CHCl₃ at 45-50°, treated simultaneously with 1.8 cc. H₂O and 5.3 cc. Br with stirring and kept 6 hrs. at 50° and 2 days at room temperature, gives 22 g. of p-bromophenylhydroxymaleimide (I), lemon-yellow, m. 239-40°; when the H₂O was omitted the yield was only 9 g.; the same yield was obtained from the Me and Bu esters. I also results from the bromination of phenylhydroxymaleimide in PhNO₂. I is stable to cold dilute KMnO₄ and to Br-H₂O. Solution of 53.6 g. I in 150 cc. H₂O containing 12.4 g. Na₂CO₃ gives the brick-red Na salt, decomp. at 321°; refluxing 5.8 g. of the salt with 2.53 g. PhCH₂Cl in 25 cc. EtOH gives 5.6 g. of the N-benzyl derivative, m. 169-70°. The brick-red gelatinous Ag salt and EtI in ether, refluxed 3 days, give the N-Et derivative, m. 191-2°. Oxidation of I gives p-BrC₆H₄CO₂H; alkaline hydrolysis of I gives NH₃, (CO₂Na)₂ and p-BrC₆H₄CH₂CO₂H. The mechanism of the formation of I is discussed.

IT 749830-37-3P

RL: SPN (Synthetic preparation); PRP (Properties); RCT (Reactant); PREP
(Preparation); RACT (Reactant or reagent)
(p-Bromophenylhydroxymaleic imide)

RN 749830-37-3 CAPLUS

CN 1H-Pyrrole-2,5-dione, 3-(4-bromophenyl)-1-hydroxy- (CA INDEX NAME)



=> log off

ALL L# QUERIES AND ANSWER SETS ARE DELETED AT LOGOFF

LOGOFF? (Y)/N/HOLD:y

STN INTERNATIONAL LOGOFF AT 13:40:31 ON 03 MAY 2009